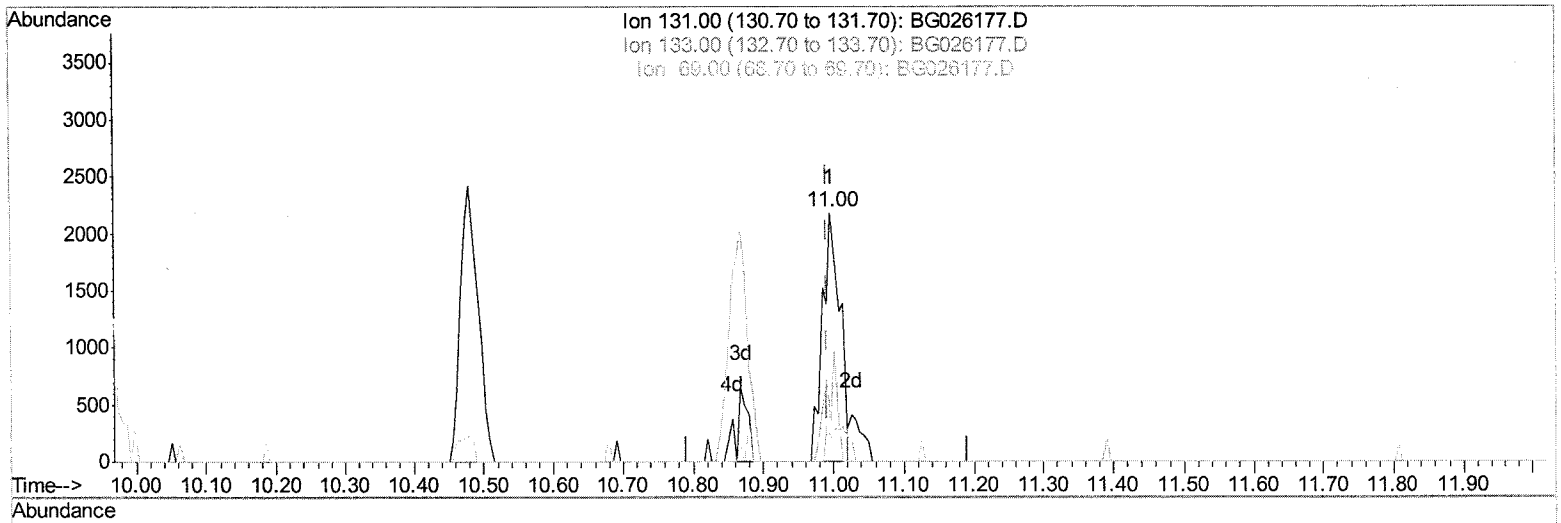
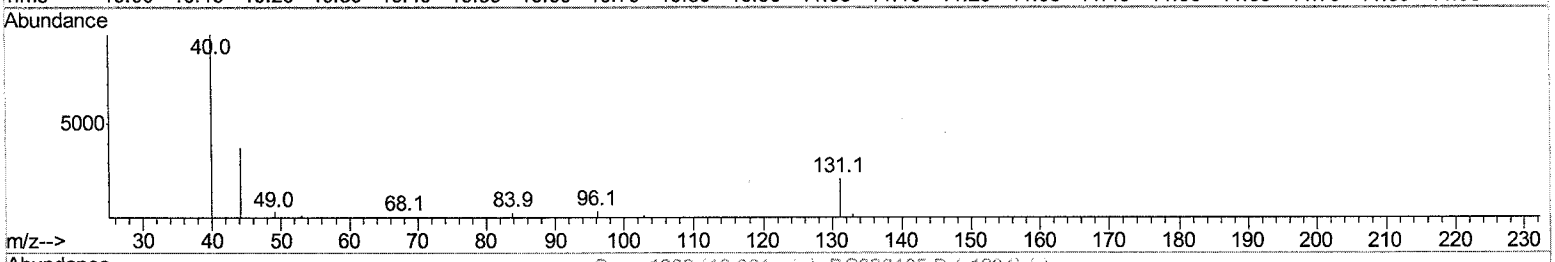
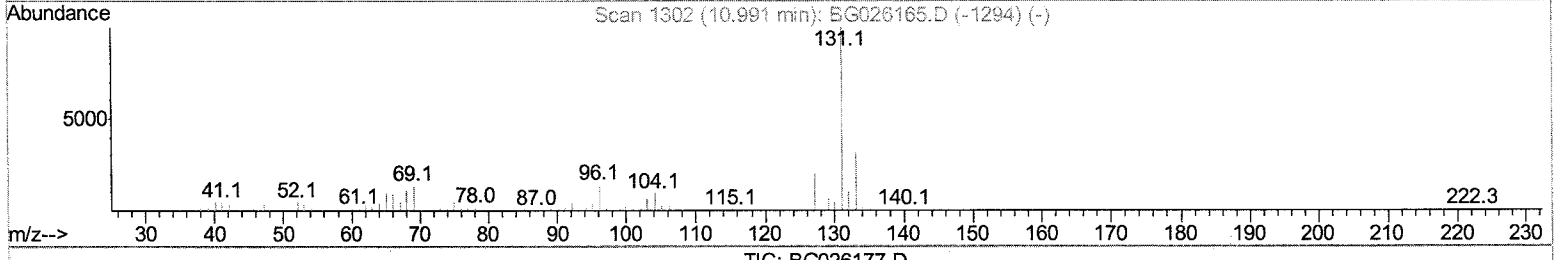
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031017\
Data File : BG026177.D
Acq On : 10 Mar 2017 17:24
Operator : SJ/MA
Sample : I2180-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 10 23:58:42 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 10 23:28:49 2017
Response via : Initial Calibration







TIC: BG026177.D

(29) 4-Chloroaniline-d4 (S)

10.997min (+0.005) 0.46ng/ul

response 3798

Ion	Exp%	
131.00	100	100
133.00	32.10	16.68#
69.00	16.20	9.83#
0.00	0.00	0.00

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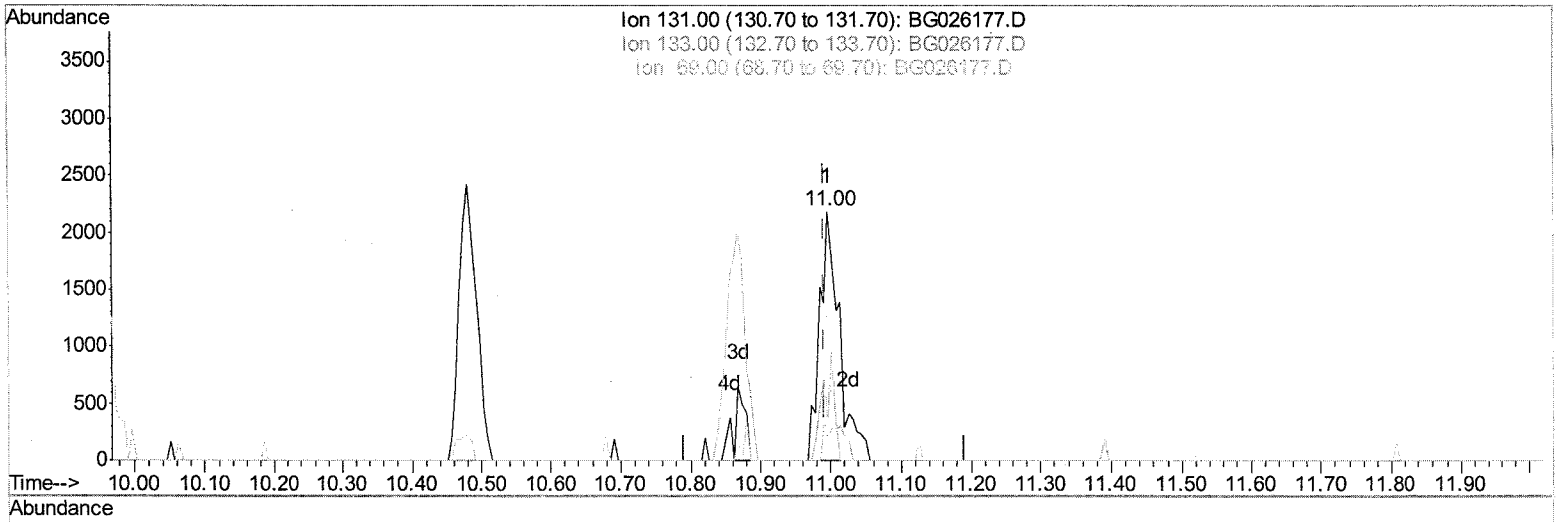

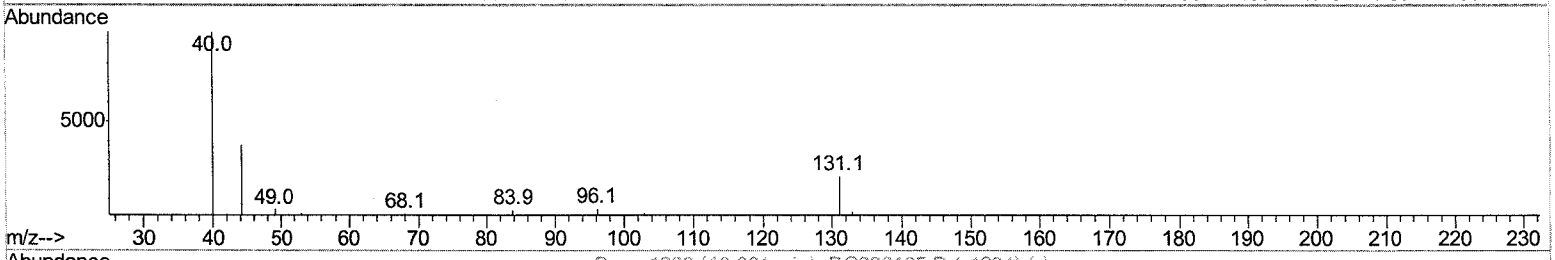

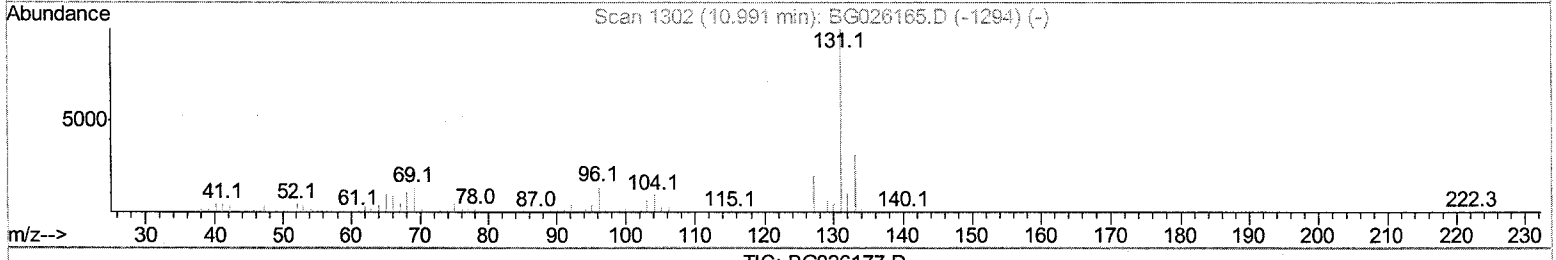

TIC: BG026177.D

(29) 4-Chloroaniline-d4 (S)

10.997min (+0.005) 0.52ng/ul m SJ 3/13/17

response 4304

Ion	Exp%	
131.00	100	100
133.00	32.10	16.68#
69.00	16.20	9.83#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_G\Data\BG031017\
Data File : BG026177.D
Acq On : 10 Mar 2017 17:24
Operator : SJ/MA
Sample : I2180-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 11 01:25:53 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 10 23:28:49 2017
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
C0BE0

Manual Integrations APPROVED
mohammad
3/13/2017 3:33:52 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	105241	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	455285	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.67	164	271368	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.40	188	664717	20.00	ng/ul	0.00
77) Chrysene-d12	21.64	240	758981	20.00	ng/ul	0.00
85) Perylene-d12	24.83	264	731605	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.49	96	7427	3.03	ng/uL	0.00
5) Phenol-d5	7.21	99	36966	4.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	126151	25.54	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	125014	19.78	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	72175	11.80	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	90183	27.42	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	76044	28.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	151624	25.11	ng/ul	0.00
29) 4-Chloroaniline-d4	11.00	131	4304m	0.52	ng/ul	0.00
43) Dimethylphthalate-d6	14.06	166	580879	32.91	ng/ul	0.00
46) Acenaphthylene-d8	14.36	160	766322	32.34	ng/ul	0.00
51) 4-Nitrophenol-d4	14.85	143	10229	2.94	ng/ul	0.00
57) Fluorene-d10	15.65	176	578256	33.22	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.76	200	62737	20.11	ng/ul	0.00
70) Anthracene-d10	17.49	188	941612	31.22	ng/ul	0.00
78) Pyrene-d10	19.77	212	1112715	35.64	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.62	264	1018406	31.58	ng/ul	0.00
Target Compounds						Qvalue
6) Phenol	7.24	94	9779	1.21	ng/ul	90

(#) = qualifier out of range (m) = manual integration (+) = signals summed

SJ 03/13/17

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Instrument :
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ClientSampleId :
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Manual Integrations APPROVED
mohammad
3/13/2017 3:33:52 PM